Amendments to the Claims:

The listing of claims will replace all prior versions, and listings, of claims in the application.

Listing of Claims:

Claim 1 (currently amended): A compound having the formula

$$R_{1}$$
 N
 Z
 R_{1}
 N
 Z
 R_{2}
 R_{3}

(1)

where:

X represents a first amine protecting group that is different from Y, wherein the first amine protecting group is selected from the group consisting of 9-fluorenylmethyl carbamate, allyl carbamate, benzyl carbamate, substituted benzyl carbamate, t-butyl carbamate, 1-adamantyl carbamate, 2-nitrobenzenesulfonyl, triphenylmethyl, (4-methoxyphenyl)diphenylmethyl, and 9-phenylfluorenyl;

Y represents a second amine protecting group that is different from X, wherein the second amine protecting group is selected from the group consisting of 9-fluorenylmethyl carbamate, allyl carbamate, benzyl carbamate, substituted benzyl carbamate, t-butyl carbamate, 1-adamantyl carbamate, 2-nitrobenzenesulfonyl, triphenylmethyl, (4-methoxyphenyl)diphenylmethyl, and 9-phenylfluorenyl;

Z represents a weak leaving group <u>selected from the group consisting of short chain alkoxides, thiolates, azide, and sulfonamides;</u>

R₁ represents an H and can be attached to the molecule at positions 2, 3 or 5;

 R_5 represents N_3 or -NH-Y;

R₆ represents a carboxylic acid; and

the stereochemical configuration at positions 2 and 4 is selected from the group consisting of (R,R), (R,S), (S,R), and (S,S).

Claims 2-3 (canceled).

Claim 4 (original): The compound of Claim 1, wherein Z is OMe.

Claim 5 (previously presented): The compound of Claim 1, wherein X is selected from the group consisting of benzyl-carbamate and t-butyl carbamate.

Claim 6 (original): The compound of Claim 1, wherein Y is 2-nitrobenzenesulfonamide. Claim 7 (original): The compound of Claim 1, wherein Y is 9-fluoroenylmethylcarbamate. Claim 8 (previously presented): The compound of Claim 1, wherein X is benzylcarbamate, R_5 is -NH-Y, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R_6 is a carboxylic acid. Claims 9-12 (canceled).

13. (withdrawn) A compound having the formula

(2)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

 R_1 represents an H, or a functional group, and can be attached to the molecule at positions 2, 3 or 5;

R₂ represents an H or a functional group;

 $\cdot R_5$ represents N_3 or NR_2X ;

 R_6 represents a carboxylic acid or a strongly activated ester; and the stereochemical configuration at positions 2 and 4 and of the carbon bearing R_1 (if R_1 is not H) can be any one of (S,S,S), (S,S,R), (S,R,S), (S,R,R), (R,S,S), (R,S,R), (R,R,S) or (R,R,R).

- 14. (withdrawn) (withdrawn) The compound of Claim 13, wherein R₅ is N₃.
- 15. (withdrawn) The compound of Claim 13, wherein R₅ is NR₂X.
- 16. (withdrawn) The compound of Claim 13, wherein Z is OMe.
- 17. (withdrawn) The compound of Claim 13, wherein X is benzylcarbamate.
- 18. (withdrawn) The compound of Claim 13, wherein Y is 2-nitrobenzenesulfonamide.
- 19. (withdrawn) The compound of Claim 13, wherein Y is 9-fluoroenylmethylcarbamate.

20. (withdrawn) The compound of Claim 13, wherein X is benzylcarbamate, R₅ is NR₂X, R₂

is H, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R_6 is a carboxylic acid.

- 21. (withdrawn) The compound of Claim 13, wherein R_1 is an alkene.
- 22. (withdrawn) The compound of Claim 13, wherein R₁ is a protected carboxylate.
- 23. (withdrawn) The compound of Claim 13, wherein R₁ is a protected alcohol.
- 24. (withdrawn) The compound of Claim 13, wherein R₁ is a protected thiol.
- 25. (withdrawn) A compound having the formula

$$R_1$$
 R_1
 R_5
 R_5
 R_6

(3)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 4 or 5;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2Y ;

 R_6 represents a carboxylic acid or a strongly activated ester; and the stereochemical configuration at positions 2 and 3 and of the carbon bearing R_1 (if R_1 is not H) can be any one of (S,S,S), (S,S,R), (S,R,S), (S,R,R), (R,S,S), (R,S,R), (R,R,S) or (R,R,R).

- 26. (withdrawn) The compound of Claim 25, wherein R₅ is N₃.
- 27. (withdrawn) The compound of Claim 25, wherein R₅ is NR₂Y.
- 28. (withdrawn) The compound of Claim 25, wherein Z is OMe.
- 29. (withdrawn) The compound of Claim 25, wherein X is benzylcarbamate.
- 30. (withdrawn) The compound of Claim 25, wherein Y is 2-nitrobenzenesulfonamide.
- 31. (withdrawn) The compound of Claim 25, wherein Y is 9-fluoroenylmethylcarbamate.

32. (withdrawn) The compound of Claim 25, wherein X is benzylcarbamate, R₅ is NR₂Y, R₂

is H, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R_6 is a carboxylic acid.

- 33. (withdrawn) The compound of Claim 25, wherein R₁ is an alkene.
- 34. (withdrawn) The compound of Claim 25, wherein R₁ is a protected carboxylate.
- 35. (withdrawn) The compound of Claim 25, wherein R₁ is a protected alcohol.
- 36. (withdrawn) The compound of Claim 25, wherein R_1 is a protected thiol.
- 37. (withdrawn) A compound having the formula

(4)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 4 or 5;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2X ;

 R_6 represents a carboxylic acid or a strongly activated ester; and the stereochemical configuration at positions 2 and 3 and of the carbon bearing R_1 (if R_1 is not H) can be any one of (S,S,S), (S,S,R), (S,R,S), (S,R,R), (R,S,S), (R,S,R), (R,R,S) or (R,R,R).

- 38. (withdrawn) The compound of Claim 37, wherein R₅ is N₃.
- 39. (withdrawn) The compound of Claim 37, wherein R₅ is NR₂X.
- 40. (withdrawn) The compound of Claim 37, wherein Z is OMe.
- 41. (withdrawn) The compound of Claim 37, wherein X is benzylcarbamate.
- 42. (withdrawn) The compound of Claim 37, wherein Y is 2-nitrobenzenesulfonamide.
- 43. (withdrawn) The compound of Claim 37, wherein Y is 9-fluoroenylmethylcarbamate.

44. (withdrawn) The compound of Claim 37, wherein X is benzylcarbamate, R₅ is NR₂X, R₂

is H, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R_6 is a carboxylic acid.

- 45. (withdrawn) The compound of Claim 37, wherein R_1 is an alkene.
- 46. (withdrawn) The compound of Claim 37, wherein R_1 is a protected carboxylate.
- 47. (withdrawn) The compound of Claim 37, wherein R₁ is a protected alcohol.
- 48. (withdrawn) The compound of Claim 37, wherein R₁ is a protected thiol.
- 49. (withdrawn) A compound having the formula

$$R_5$$
 Q
 Z
 R_1
 Q
 Z
 R_6

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 3, 5 or 6;

(5)

R₂ represents an H or a functional group;

R₅ represents N₃ or NR₂Y;

 R_6 represents a carboxylic acid or a strongly activated ester; and the stereochemical configuration at positions 2 and 4 and of the carbon bearing R_1 (if R_1 is not H) can be any one of (S,S,S), (S,S,R), (S,R,S), (S,R,R), (R,S,S), (R,S,R), (R,S,R), or (R,R,R).

- 50. (withdrawn) The compound of Claim 49, wherein R₅ is N₃.
- 51. (withdrawn) The compound of Claim 49, wherein R₅ is NR₂Y.
- 52. (withdrawn) The compound of Claim 49, wherein Z is OMe.
- 53. (withdrawn) The compound of Claim 49, wherein X is benzylcarbamate.
- 54. (withdrawn) The compound of Claim 49, wherein Y is 2-nitrobenzenesulfonamide.
- 55. (withdrawn) The compound of Claim 49, wherein Y is 9-fluoroenylmethylcarbamate.

56. (withdrawn) The compound of Claim 49, wherein X is benzylcarbamate, R₅ is NR₂Y, R₂

is H, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R_6 is a carboxylic acid.

- 57. (withdrawn) The compound of Claim 49, wherein R₁ is an alkene.
- 58. (withdrawn) The compound of Claim 49, wherein R₁ is a protected carboxylate.
- 59. (withdrawn) The compound of Claim 49, wherein R_1 is a protected alcohol.
- 60. (withdrawn) The compound of Claim 49, wherein R₁ is a protected thiol.
- 61. (withdrawn) A compound having the formula

$$R_1$$
 R_1
 R_1
 R_1
 R_1
 R_2
 R_3
 R_4
 R_5
 R_6
 R_6
 R_7
 R_6

(6)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 3, 5 or 6;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2X ;

 $R_{\rm 6}$ represents a carboxylic acid or a strongly activated ester ; and

the stereochemical configuration at positions 2 and 4 and of the carbon bearing R_1 (if R_1 is not H) can be any one of (S,S,S), (S,S,R), (S,R,S), (S,R,R), (R,S,S), (R,S,R), (R,R,S) or (R,R,R).

- 62. (withdrawn) The compound of Claim 61, wherein R₅ is N₃.
- 63. (withdrawn) The compound of Claim 61, wherein R₅ is NR₂X.
- 64. (withdrawn) The compound of Claim 61, wherein Z is OMe.
- 65. (withdrawn) The compound of Claim 61, wherein X is benzylcarbamate.
- 66. (withdrawn) The compound of Claim 61, wherein Y is 2-nitrobenzenesulfonamide.
- 67. (withdrawn) The compound of Claim 61, wherein Y is 9-fluoroenylmethylcarbamate.

68. (withdrawn) The compound of Claim 61, wherein X is benzylcarbamate, R₅ is NR₂X, R₂

is H, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R_6 is a carboxylic acid.

- 69. (withdrawn) The compound of Claim 61, wherein R_1 is an alkene.
- 70. (withdrawn) The compound of Claim 61, wherein R₁ is a protected carboxylate.
- 71. (withdrawn) The compound of Claim 61, wherein R₁ is a protected alcohol.
- 72. (withdrawn) The compound of Claim 61, wherein R_1 is a protected thiol.
- 73. (withdrawn) A compound having the formula

(7)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 3, 4 or 6;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2Y ;

R₆ represents a carboxylic acid or a strongly activated ester; and

the stereochemical configuration at positions 2 and 5 and of the carbon bearing R_1 (if R_1 is not H) can be any one of (S,S,S), (S,S,R), (S,R,S), (S,R,R), (R,S,S), (R,S,R), (R,R,S) or (R,R,R).

- 74. (withdrawn) The compound of Claim 73, wherein R_5 is N_3 .
- 75. (withdrawn) The compound of Claim 73, wherein R_5 is NR_2Y .
- 76. (withdrawn) The compound of Claim 73, wherein Z is OMe.
- 77. (withdrawn) The compound of Claim 73, wherein X is benzylcarbamate.
- 78. (withdrawn) The compound of Claim 73, wherein Y is 2-nitrobenzenesulfonamide.
- 79. (withdrawn) The compound of Claim 73, wherein Y is 9-fluoroenylmethylcarbamate.

80. (withdrawn) The compound of Claim 73, wherein X is benzylcarbamate, R₅ is NR₂Y, R₂

is H, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R_6 is a carboxylic acid.

81. (withdrawn) The compound of Claim 73, wherein R_1 is an alkene.

82. (withdrawn) The compound of Claim 73, wherein R₁ is a protected carboxylate.

83. (withdrawn) The compound of Claim 73, wherein R_1 is a protected alcohol.

84. (withdrawn) The compound of Claim 73, wherein R₁ is a protected thiol.

85. (withdrawn) A compound having the formula

$$\begin{array}{c} R_6 \\ R_5 \\ \hline \\ N \\ V \\ O \end{array}$$

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 3, 5 or 6;

(8)

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2X ;

R₆ represents a carboxylic acid or a strongly activated ester; and

the stereochemical configuration at positions 2 and 5 and of the carbon bearing R_1 (if R_1 is not H) can be any one of (S,S,S), (S,S,R), (S,R,S), (S,R,R), (R,S,S), (R,S,R), (R,R,S) or (R,R,R).

86. (withdrawn) The compound of Claim 85, wherein R₅ is N₃.

87. (withdrawn) The compound of Claim 85, wherein R₅ is NR₂X.

88. (withdrawn) The compound of Claim 85, wherein Z is OMe.

89. (withdrawn) The compound of Claim 85, wherein X is benzylcarbamate.

90. (withdrawn) The compound of Claim 85, wherein Y is 2-nitrobenzenesulfonamide.

91. (withdrawn) The compound of Claim 85, wherein Y is 9-fluoroenylmethylcarbamate.

92. (withdrawn) The compound of Claim 85, wherein X is benzylcarbamate, R₅ is NR₂X, R₂

is H, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R₆ is a carboxylic acid.

93. (withdrawn) The compound of Claim 85, wherein R₁ is an alkene.

94. (withdrawn) The compound of Claim 85, wherein R₁ is a protected carboxylate.

95. (withdrawn) The compound of Claim 85, wherein R_1 is a protected alcohol.

96. (withdrawn) The compound of Claim 85, wherein R_1 is a protected thiol.

97. (withdrawn) A compound having the formula

(9)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

 R_1 represents an H, or a functional group, and can be attached to the molecule at positions 2, 4, 5 or 6;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2Y ;

R₆ represents a carboxylic acid or a strongly activated ester; and

the stereochemical configuration at positions 2 and 3 and of the carbon bearing R_1 (if R_1 is not H) can be any one of (S,S,S), (S,S,R), (S,R,S), (S,R,R), (R,S,S), (R,S,R), (R,R,S) or (R,R,R).

98. (withdrawn) The compound of Claim 97, wherein R₅ is N₃.

99. (withdrawn) The compound of Claim 97, wherein R₅ is NR₂Y.

100. (withdrawn) The compound of Claim 97, wherein Z is OMe.

101. (withdrawn) The compound of Claim 97, wherein X is benzylcarbamate.

102. (withdrawn) The compound of Claim 97, wherein Y is 2-nitrobenzenesulfonamide.

103. (withdrawn) The compound of Claim 97, wherein Y is 9-fluoroenylmethylcarbamate.

104. (withdrawn) The compound of Claim 97, wherein X is benzylcarbamate, R₅ is NR₂Y,

 R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is –OMe, and R_6 is a carboxylic acid.

105. (withdrawn) The compound of Claim 97, wherein R_1 is an alkene.

106. (withdrawn) The compound of Claim 97; wherein R_1 is a protected carboxylate.

107. (withdrawn) The compound of Claim 97, wherein R₁ is a protected alcohol.

108. (withdrawn) The compound of Claim 97, wherein R₁ is a protected thiol.

109. (withdrawn) A compound having the formula

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

(10)

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 4, 5 or 6;

R₂ represents an H or a functional group;

R₅ represents N₃ or NR₂X;

 R_6 represents a carboxylic acid or a strongly activated ester; and the stereochemical configuration at positions 2 and 3 and of the carbon bearing R_1 (if R_1 is not H) can be any one of (S,S,S), (S,S,R), (S,R,S), (S,R,R), (R,S,S), (R,S,R), (R,R,S) or (R,R,R).

110. (withdrawn) The compound of Claim 109, wherein R₅ is N₃.

111. (withdrawn) The compound of Claim 109, wherein R₅ is NR₂X.

112. (withdrawn) The compound of Claim 109, wherein Z is OMe.

113. (withdrawn) The compound of Claim 109, wherein X is benzylcarbamate.

114. (withdrawn) The compound of Claim 109, wherein Y is 2-nitrobenzenesulfonamide.

115. (withdrawn) The compound of Claim 109, wherein Y is 9-fluoroenylmethylcarbamate.

116. (withdrawn) The compound of Claim 109, wherein X is benzylcarbamate, R₅ is NR₂X,

 R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is –OMe, and R_6 is a carboxylic acid.

117. (withdrawn) The compound of Claim 109, wherein R₁ is an alkene.

118. (withdrawn) The compound of Claim 109, wherein R₁ is a protected carboxylate.

119. (withdrawn) The compound of Claim 109, wherein R_1 is a protected alcohol.

120. (withdrawn) The compound of Claim 109, wherein R₁ is a protected thiol.

121. (withdrawn) A compound having the formula

$$\begin{array}{c|c}
R_1 & 4 \\
\hline
 & 7 \\
O & Z
\end{array}$$

$$\begin{array}{c}
R_1 & 4 \\
\hline
 & 2 \\
\hline
 & X
\end{array}$$

(11)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions

2, 3, 4, 5, 6, 8 or 9;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2Y ;

R₆ represents a carboxylic acid or a strongly activated ester; and

the stereochemical configuration at positions 2, 4, 7, 9 and of the carbon bearing R_1 (if R_1 is not H) can be any of the 32 combinations of (R) and (S).

122. (withdrawn) The compound of Claim 121, wherein R₅ is N₃.

123. (withdrawn) The compound of Claim 121, wherein R₅ is NR₂Y.

124. (withdrawn) The compound of Claim 121, wherein Z is OMe.

125. (withdrawn) The compound of Claim 121, wherein X is benzylcarbamate.

126. (withdrawn) The compound of Claim 121, wherein Y is 2-nitrobenzenesulfonamide.

127. (withdrawn) The compound of Claim 121, wherein Y is 9-fluoroenylmethylcarbamate.

128. (withdrawn) The compound of Claim 121, wherein X is benzylcarbamate, R₅ is NR₂Y,

 R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is –OMe, and R_6 is a carboxylic acid.

129. (withdrawn) The compound of Claim 121, wherein R_1 is an alkene.

130. (withdrawn) The compound of Claim 121, wherein R_1 is a protected carboxylate.

131. (withdrawn) The compound of Claim 121, wherein R_1 is a protected alcohol.

132. (withdrawn) The compound of Claim 121, wherein R_1 is a protected thiol.

$$R_5$$
 R_6 R_6

(12)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 3, 4, 5, 6, 8 or 9;

R₂ represents an H or a functional group;

R₅ represents N₃ or NR₂X;

 R_6 represents a carboxylic acid or a strongly activated ester; and the stereochemical configuration at positions 2, 4, 7, 9 and of the carbon bearing R_1 (if R_1 is not H) can be any of the 32 combinations of (R) and (S).

- 134. (withdrawn) The compound of Claim 133, wherein R₅ is N₃.
- 135. (withdrawn) The compound of Claim 133, wherein R₅ is NR₂X.
- 136. (withdrawn) The compound of Claim 133, wherein Z is OMe.
- 137. (withdrawn) The compound of Claim 133, wherein X is benzylcarbamate.
- 138. (withdrawn) The compound of Claim 133, wherein Y is 2-nitrobenzenesulfonamide.
- 139. (withdrawn) The compound of Claim 133, wherein Y is 9-fluoroenylmethylcarbamate.
- 140. (withdrawn) The compound of Claim 133, wherein X is benzylcarbamate, R₅ is NR₂X,

 R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R_6 is a carboxylic acid.

- 141. (withdrawn) The compound of Claim 133, wherein R₁ is an alkene.
- 142. (withdrawn) The compound of Claim 133, wherein R₁ is a protected carboxylate.
- 143. (withdrawn) The compound of Claim 133, wherein R₁ is a protected alcohol.
- 144. (withdrawn) The compound of Claim 133, wherein R_1 is a protected thiol.

(13)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions

2, 3, 5, 6, 7 or 8;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2Y ;

R₆ represents a carboxylic acid or a strongly activated ester; and

the stereochemical configuration at the positions 2, 3, 4 and 5, and of the carbon bearing

 R_1 (if R_1 is not H) can be any of the 32 combinations of (R) and (S).

146. (withdrawn) The compound of Claim 145, wherein R₅ is N₃.

147. (withdrawn) The compound of Claim 145, wherein R₅ is NR₂Y.

148. (withdrawn) The compound of Claim 145, wherein Z is OMe.

149. (withdrawn) The compound of Claim 145, wherein X is benzylcarbamate.

150. (withdrawn) The compound of Claim 145, wherein Y is 2-nitrobenzenesulfonamide.

151. (withdrawn) The compound of Claim 145, wherein Y is 9-fluoroenylmethylcarbamate.

152. (withdrawn) The compound of Claim 145, wherein X is benzylcarbamate, R₅ is NR₂Y,

 R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R_6 is a carboxylic acid.

153. (withdrawn) The compound of Claim 145, wherein R₁ is an alkene.

154. (withdrawn) The compound of Claim 145, wherein R₁ is a protected carboxylate.

155. (withdrawn) The compound of Claim 145, wherein R₁ is a protected alcohol.

156. (withdrawn) The compound of Claim 145, wherein R₁ is a protected thiol.

$$\begin{array}{c|c}
R_6 & 4 \\
0 & 3 \\
2 & 8 \\
7 & 7
\end{array}$$

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 3, 5, 6, 7 or 8;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2X ;

R₆ represents a carboxylic acid or a strongly activated ester; and

(14)

the stereochemical configuration at the positions 2, 3, 4 and 5, and of the carbon bearing R_1 (if R_1 is not H) can be any of the 32 combinations of (R) and (S).

158. (withdrawn) The compound of Claim 157, wherein R₅ is N₃.

159. (withdrawn) The compound of Claim 157, wherein R₅ is NR₂X.

160. (withdrawn) The compound of Claim 157, wherein Z is OMe.

161. (withdrawn) The compound of Claim 157, wherein X is benzylcarbamate.

162. (withdrawn) The compound of Claim 157, wherein Y is 2-nitrobenzenesulfonamide.

163. (withdrawn) The compound of Claim 157, wherein Y is 9-fluoroenylmethylcarbamate.

164. (withdrawn) The compound of Claim 157, wherein X is benzylcarbamate, R₅ is NR₂X,

 R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is –OMe, and R_6 is a carboxylic acid.

165. (withdrawn) The compound of Claim 157, wherein R_1 is an alkene.

166. (withdrawn) The compound of Claim 157, wherein R_1 is a protected carboxylate.

167. (withdrawn) The compound of Claim 157, wherein R_1 is a protected alcohol.

168. (withdrawn) The compound of Claim 157, wherein R_1 is a protected thiol.

169. (withdrawn) A method of synthesizing bis peptides comprising the steps of:

1) providing a solid support;

- 2) activating a first bis amino acid or naturally occurring amino acid;
- 3) attaching the bis amino acid or naturally occurring amino acid to the support;
- 4) removing the leading edge amine protecting group if a *bis* amino acid is used, or the amine protecting group if a naturally occurring amino acid is used;
- 5) activating and attaching a next bis amino acid or a next naturally occurring amino acid to the leading edge amine of the bis amino acid or amine of the naturally occurring amino acid; and
 - 6) repeating steps 4 and 5 as necessary to achieve the desired chain length;
 - 7) detaching the synthesized bis peptide from the support; and
 - 8) isolating the synthesized bis peptide,

where the *bis* peptide synthesized in the above manner has at least two contiguous *bis* amino acids, and a rigidification step is carried out either after step 4 or after detachment of the *bis* peptide from the solid support.

170. (withdrawn) The method of Claim 169, further comprising the step of modifying or adding a functional group, after step 5.

- 171. (withdrawn) A method of synthesizing bis peptides comprising the steps of:
- 1) providing a *bis*-amino acid or *bis*-peptide fragment containing a mixture of *bis*-amino acid and naturally occurring amino acid with an unprotected leading edge amine and a protected trailing edge carboxylic acid;
- 2) providing a *bis*-s or *bis*-peptide fragment containing a mixture of *bis*-amino acid and naturally occurring amino acids with a protected leading edge amine and an activated ester;
 - 3) coupling the two fragments in solution;
 - 4) isolating the synthesized bis-peptide;
- 5) removing the leading edge amine protecting group or the trailing end carboxylic acid protecting group; and
 - 6) repeating steps 1,2,3,4 to achieve the desired chain length;

where the *bis* peptide synthesized in the above manner has at least two contiguous *bis* amino acids, and a rigidification step is carried out either after step 3 or after detachment of the *bis* peptide from the solid support.

172. (withdrawn) The method of Claim 171, further comprising the step of modifying or adding a functional group, after step 3.

$$\begin{array}{c|c}
Z & O \\
 & 4 & R_5 \\
 & R_6 & 9 & R_1 \\
 & & 7 & \\
 & & X & (15)
\end{array}$$

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

 R_1 represents an H, or a functional group, and can be attached to the molecule at positions 2, 3, 5, 6, 7, 8 or 9;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2Y ;

 $R_{\rm 6}$ represents a carboxylic acid or a strongly activated ester ; and

the stereochemical configuration at the positions 2, 3, 4 and 5, and of the carbon bearing

 R_1 (if R_1 is not H) can be any of the 32 combinations of (R) and (S).

174. (withdrawn) The compound of Claim 173, wherein R₅ is N₃.

175. (withdrawn) The compound of Claim 173, wherein R₅ is NR₂Y.

176. (withdrawn) The compound of Claim 173, wherein Z is OMe.

177. (withdrawn) The compound of Claim 173, wherein X is benzylcarbamate.

178. (withdrawn) The compound of Claim 173, wherein Y is 2-nitrobenzenesulfonamide.

179. (withdrawn) The compound of Claim 173, wherein Y is 9-fluoroenylmethylcarbamate.

180. (withdrawn) The compound of Claim 173, wherein X is benzylcarbamate, R₅ is NR₂Y,

 R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is –OMe, and R_6 is a carboxylic acid.

181. (withdrawn) The compound of Claim 173, wherein R_1 is an alkene.

182. (withdrawn) The compound of Claim 173, wherein R_1 is a protected carboxylate.

183. (withdrawn) The compound of Claim 173, wherein R_1 is a protected alcohol.

184. (withdrawn) The compound of Claim 173, wherein R_1 is a protected thiol.

$$\begin{array}{c|c}
R_6 & 4 \\
\hline
0 & 3 \\
\hline
2 & 3 \\
\hline
9 \\
R_1 \\
7 \\
\hline
7
\end{array}$$

(16)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 3, 5, 6, 7, 8 or 9;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2X ;

 R_6 represents a carboxylic acid or a strongly activated ester; and

the stereochemical configuration at the positions 2, 3, 4 and 5, and of the carbon bearing

 R_1 (if R_1 is not H) can be any of the 32 combinations of (R) and (S).

186. (withdrawn) The compound of Claim 185, wherein R_5 is N_3 .

187. (withdrawn) The compound of Claim 185, wherein R₅ is NR₂X.

188. (withdrawn) The compound of Claim 185, wherein Z is OMe.

189. (withdrawn) The compound of Claim 185, wherein X is benzylcarbamate.

190. (withdrawn) The compound of Claim 185, wherein Y is 2-nitrobenzenesulfonamide.

191. (withdrawn) The compound of Claim 185, wherein Y is 9-fluoroenylmethylcarbamate.

192. (withdrawn) The compound of Claim 185, wherein X is benzylcarbamate, R₅ is NR₂X,

 R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R_6 is a carboxylic acid.

193. (withdrawn) The compound of Claim 185, wherein R₁ is an alkene.

194. (withdrawn) The compound of Claim 185, wherein R₁ is a protected carboxylate.

195. (withdrawn) The compound of Claim 185, wherein R_1 is a protected alcohol.

196. (withdrawn) The compound of Claim 185, wherein R₁ is a protected thiol.

(17)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 3, 4, 6, 7, 8 or 9;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2Y ;

R₆ represents a carboxylic acid or a strongly activated ester; and

the stereochemical configuration at the positions 2, 4, 5 and 6, and of the carbon bearing

 R_1 (if R_1 is not H) can be any of the 32 combinations of (R) and (S).

198. (withdrawn) The compound of Claim 197, wherein R₅ is N₃.

199. (withdrawn) The compound of Claim 197, wherein R₅ is NR₂Y.

200. (withdrawn) The compound of Claim 197, wherein Z is OMe.

201. (withdrawn) The compound of Claim 197, wherein X is benzylcarbamate.

202. (withdrawn) The compound of Claim 197, wherein Y is 2-nitrobenzenesulfonamide.

203. (withdrawn) The compound of Claim 197, wherein Y is 9-fluoroenylmethylcarbamate.

204. (withdrawn) The compound of Claim 197, wherein X is benzylcarbamate, R₅ is NR₂Y,

 R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is –OMe, and R_6 is a carboxylic acid.

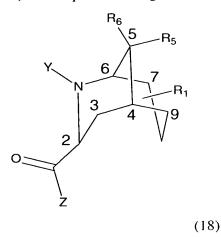
205. (withdrawn) The compound of Claim 197, wherein R₁ is an alkene.

206. (withdrawn) The compound of Claim 197, wherein R₁ is a protected carboxylate.

207. (withdrawn) The compound of Claim 197, wherein R₁ is a protected alcohol.

208. (withdrawn) The compound of Claim 197, wherein R₁ is a protected thiol.

209. (withdrawn) A compound having the formula



where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 3, 4, 6, 7, 8 or 9;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2X ;

R₆ represents a carboxylic acid or a strongly activated ester; and

the stereochemical configuration at the positions 2, 4, 5 and 6, and of the carbon bearing R_1 (if R_1 is not H) can be any of the 32 combinations of (R) and (S).

210. (withdrawn) The compound of Claim 209, wherein R₅ is N₃.

211. (withdrawn) The compound of Claim 209, wherein R₅ is NR₂X.

212. (withdrawn) The compound of Claim 209, wherein Z is OMe.

213. (withdrawn) The compound of Claim 209, wherein X is benzylcarbamate.

214. (withdrawn) The compound of Claim 209, wherein Y is 2-nitrobenzenesulfonamide.

215. (withdrawn) The compound of Claim 209, wherein Y is 9-fluoroenylmethylcarbamate.

216. (withdrawn) The compound of Claim 209, wherein X is benzylcarbamate, R₅ is NR₂X,

 R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R_6 is a carboxylic acid.

217. (withdrawn) The compound of Claim 209, wherein R₁ is an alkene.

218. (withdrawn) The compound of Claim 209, wherein R₁ is a protected carboxylate.

- 219. (withdrawn) The compound of Claim 209, wherein R₁ is a protected alcohol.
- 220. (withdrawn) The compound of Claim 209, wherein R₁ is a protected thiol.
- 221. (withdrawn) A compound having the formula

$$R_6$$
 R_5
 R_5
 R_5
 R_7
 R_7

(19)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 3, 4, 5, 7 or 8;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2Y ;

R₆ represents a carboxylic acid or a strongly activated ester; and

the stereochemical configuration at the positions 2, 4, 6 and 8, and of the carbon bearing R_1 (if R_1 is not H) can be any of the 32 combinations of (R) and (S).

- 222. (withdrawn) The compound of Claim 221, wherein R₅ is N₃.
- 223. (withdrawn) The compound of Claim 221, wherein R₅ is NR₂Y.
- 224. (withdrawn) The compound of Claim 221, wherein Z is OMe.
- 225. (withdrawn) The compound of Claim 221, wherein X is benzylcarbamate.
- 226. (withdrawn) The compound of Claim 221, wherein Y is 2-nitrobenzenesulfonamide.
- 227. (withdrawn) The compound of Claim 221, wherein Y is 9-fluoroenylmethylcarbamate.
- 228. (withdrawn) The compound of Claim 221, wherein X is benzylcarbamate, R₅ is NR₂Y,
- R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R_6 is a carboxylic acid.
- 229. (withdrawn) The compound of Claim 221, wherein R₁ is an alkene.
- 230. (withdrawn) The compound of Claim 221, wherein R₁ is a protected carboxylate.
- 231. (withdrawn) The compound of Claim 221, wherein R₁ is a protected alcohol.

232. (withdrawn) The compound of Claim 221, wherein R₁ is a protected thiol.

233. (withdrawn) A compound having the formula

(20)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions

2, 3, 4, 5, 7 or 8;

R₂ represents an H or a functional group;

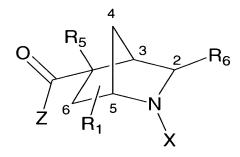
 R_5 represents N_3 or NR_2X ;

R₆ represents a carboxylic acid or a strongly activated ester; and

the stereochemical configuration at the positions 2, 4, 6 and 8, and of the carbon bearing

 R_1 (if R_1 is not H) can be any of the 32 combinations of (R) and (S).

- 234. (withdrawn) The compound of Claim 233, wherein R₅ is N₃.
- 235. (withdrawn) The compound of Claim 233, wherein R₅ is NR₂X.
- 236. (withdrawn) The compound of Claim 233, wherein Z is OMe.
- 237. (withdrawn) The compound of Claim 233, wherein X is benzylcarbamate.
- 238. (withdrawn) The compound of Claim 233, wherein Y is 2-nitrobenzenesulfonamide.
- 239. (withdrawn) The compound of Claim 233, wherein Y is 9-fluoroenylmethylcarbamate.
- 240, (withdrawn) The compound of Claim 233, wherein X is benzylcarbamate, R₅ is NR₂X,
- R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R_6 is a carboxylic acid.
- 241. (withdrawn) The compound of Claim 233, wherein R_1 is an alkene.
- 242. (withdrawn) The compound of Claim 233, wherein R_1 is a protected carboxylate.
- 243. (withdrawn) The compound of Claim 233, wherein R₁ is a protected alcohol.
- 244. (withdrawn) The compound of Claim 233, wherein R₁ is a protected thiol.



(21)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 3, 4, 5 or 6;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2Y ;

 R_6 represents a carboxylic acid or a strongly activated ester; and the stereochemical configuration at the positions 2, 3, 5 and 7, and of the carbon bearing

 R_1 (if R_1 is not H) can be any of the 32 combinations of (R) and (S).

246. (withdrawn) The compound of Claim 245, wherein R_5 is N_3 .

247. (withdrawn) The compound of Claim 245, wherein R₅ is NR₂Y.

248. (withdrawn) The compound of Claim 245, wherein Z is OMe.

249. (withdrawn) The compound of Claim 245, wherein X is benzylcarbamate.

250. (withdrawn) The compound of Claim 245, wherein Y is 2-nitrobenzenesulfonamide.

251. (withdrawn) The compound of Claim 245, wherein Y is 9-fluoroenylmethylcarbamate.

252. (withdrawn) The compound of Claim 245, wherein X is benzylcarbamate, R₅ is NR₂Y,

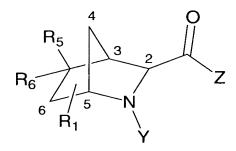
 R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is –OMe, and R_6 is a carboxylic acid.

253. (withdrawn) The compound of Claim 245, wherein R₁ is an alkene.

254. (withdrawn) The compound of Claim 245, wherein R₁ is a protected carboxylate.

255. (withdrawn) The compound of Claim 245, wherein R₁ is a protected alcohol.

256. (withdrawn) The compound of Claim 245, wherein R₁ is a protected thiol.



(22)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 3, 4, 5 or 6;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2X ;

R₆ represents a carboxylic acid or a strongly activated ester; and

the stereochemical configuration at the positions 2, 3, 5 and 7, and of the carbon bearing

 R_1 (if R_1 is not H) can be any of the 32 combinations of (R) and (S).

258. (withdrawn) The compound of Claim 257, wherein R_5 is N_3 .

259. (withdrawn) The compound of Claim 257, wherein R₅ is NR₂X.

260. (withdrawn) The compound of Claim 257, wherein Z is OMe.

261. (withdrawn) The compound of Claim 257, wherein X is benzylcarbamate.

262. (withdrawn) The compound of Claim 257, wherein Y is 2-nitrobenzenesulfonamide.

263. (withdrawn) The compound of Claim 257, wherein Y is 9-fluoroenylmethylcarbamate.

264. (withdrawn) The compound of Claim 257, wherein X is benzylcarbamate, R₅ is NR₂X,

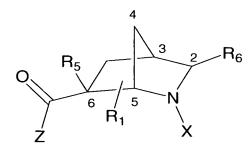
 R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R_6 is a carboxylic acid.

265. (withdrawn) The compound of Claim 257, wherein R₁ is an alkene.

266. (withdrawn) The compound of Claim 257, wherein R₁ is a protected carboxylate.

267. (withdrawn) The compound of Claim 257, wherein R₁ is a protected alcohol.

268. (withdrawn) The compound of Claim 257, wherein R₁ is a protected thiol.



(23)

where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

R₁ represents an H, or a functional group, and can be attached to the molecule at positions 2, 3, 4, 5 or 7;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2Y ;

 R_6 represents a carboxylic acid or a strongly activated ester; and the stereochemical configuration at the positions 2, 3, 5 and 6, and of the carbon bearing R_1 (if R_1 is not H) can be any of the 32 combinations of (R) and (S).

270. (withdrawn) The compound of Claim 269, wherein R₅ is N₃.

271. (withdrawn) The compound of Claim 269, wherein R₅ is NR₂Y.

272. (withdrawn) The compound of Claim 269, wherein Z is OMe.

273. (withdrawn) The compound of Claim 269, wherein X is benzylcarbamate.

274. (withdrawn) The compound of Claim 269, wherein Y is 2-nitrobenzenesulfonamide.

275. (withdrawn) The compound of Claim 269, wherein Y is 9-fluoroenylmethylcarbamate.

276. (withdrawn) The compound of Claim 269, wherein X is benzylcarbamate, R₅ is NR₂Y,

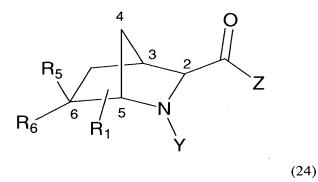
 R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is –OMe, and R_6 is a carboxylic acid.

277. (withdrawn) The compound of Claim 269, wherein R₁ is an alkene.

278. (withdrawn) The compound of Claim 269, wherein R₁ is a protected carboxylate.

279. (withdrawn) The compound of Claim 269, wherein R₁ is a protected alcohol.

280. (withdrawn) The compound of Claim 269, wherein R_1 is a protected thiol.



where:

X represents a first amine protecting group;

Y represents a second amine protecting group;

Z represents a weak leaving group;

 R_1 represents an H, or a functional group, and can be attached to the molecule at positions 2, 3, 4, 5 or 7;

R₂ represents an H or a functional group;

 R_5 represents N_3 or NR_2X ;

 R_6 represents a carboxylic acid or a strongly activated ester; and the stereochemical configuration at the positions 2, 3, 5 and 6, and of the carbon bearing R_1 (if R_1 is not H) can be any of the 32 combinations of (R) and (S).

- 282. (withdrawn) The compound of Claim 281, wherein R_5 is N_3 .
- 283. (withdrawn) The compound of Claim 281, wherein R₅ is NR₂X.
- 284. (withdrawn) The compound of Claim 281, wherein Z is OMe.
- 285. (withdrawn) The compound of Claim 281, wherein X is benzylcarbamate.
- 286. (withdrawn) The compound of Claim 281, wherein Y is 2-nitrobenzenesulfonamide.
- 287. (withdrawn) The compound of Claim 281, wherein Y is 9-fluoroenylmethylcarbamate.
- 288. (withdrawn) The compound of Claim 281, wherein X is benzylcarbamate, R₅ is NR₂X,
- R_2 is H, Y is 9-fluoroenylmethylcarbamate, Z is -OMe, and R_6 is a carboxylic acid.
- 289. (withdrawn) The compound of Claim 281, wherein R_1 is an alkene.
- 290. (withdrawn) The compound of Claim 281, wherein R₁ is a protected carboxylate.
- 291. (withdrawn) The compound of Claim 281, wherein R₁ is a protected alcohol.
- 292. (withdrawn) The compound of Claim 281, wherein R₁ is a protected thiol.